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Standard Form 298 (Rev. 8-98) Prescribed by ANSI Std. 239.18

MEMORANDUM FOR PR (In-House Publication)

FROM: PROI (TI) (STINFO)

18 May 2000

SUBJECT: Authorization for Release of Technical Information, Control Number: AFRL-PR-ED-TP-2000-117 Perkins, Dr. Leslie, "Computational Chemistry and Materials Science CTA"

High Performance Computing Mod. Office (no location given, 27 Mar 00)	(Statement A) (Submission Deadline: N/A)
 b.) military/national critical technology, c.) export cond.) appropriateness for release to a foreign nation, and Comments: 	e.) technical sensitivity and/or economic sensitivity.
Signature	
This request has been reviewed by the Public Affair and/or b) possible higher headquarters review. Comments:	
Signature	Date
e.) parallel review completed if required, and f.) formated Comments:	ary/national critical technology, d.) economic sensitivity, at and completion of meeting clearance form if required
Signature	Date
4. This request has been reviewed by PR for: a.) techn appropriateness of distribution statement, d.) technical national critical technology, and f.) data rights and pat Comments:	I sensitivity and economic sensitivity, e.) military/ sentability
	APPROVED/APPROVED AS AMENDED/DISAPPROVED

LESLIE. S. PERKINS, Ph.D **Staff Scientist**

Propulsion Directorate

(Date)

PR2000 CHSSI Review

Computational Chemistry and Materials Science CTA

Dr. Leslie S. Perkins CTA Leader U.S. Air Force Research Laboratory Propulsion Directorate Edwards AFB, CA

 $oldsymbol{CM}_{ ext{Computational Chemistry}}$ and Materials Science

Distribution A: Approved for public release distribution unlimited.

CCM Concerns within DoD

- Conductive and insulative materials for communications systems and sensors
- Aging and Surveillance of DoD stockpiles
- Contaminant decomposition within a known environment
- Energetic fuels and oxidizers for propulsion in critical military systems
- Low-observable coatings to protect the warfighter

Project Summary

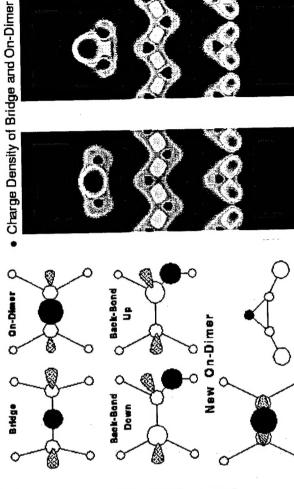
	Scalability		
Project	Range	Platforms	Deliverables
CCM-1:	23% - 70%	Origin 2000	ACRES
Car-Parrinello Methods for Solids		IBM SP	DoDPW
CCM-2:	%08 - %09	IBM SP	GAMESS
Quantum Chemistry		Cray T3E	SAPT
CCM-3:	62% - 97%	Origin 2000	TBLibrary
Tight Binding Molecular		IBM SP	Static
Dynamics			OWIG I
CCM-4:	44% - 80%	Origin 2000	FMD
Classical Molecular Dynamics		Cray T3E	PIMD/CMD

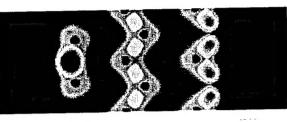
 $oldsymbol{CM}_{ ext{Computational Chemistry}}$ and Materials Science

CCM-1: Car-Parrinello Methods for Solids

Synopsis

- codes (Car-Parrinello) to be used in DoD • Objective: Produce density functional materials research
- Partners: Harvard, NRL
- Development Paradigm:
- DoDPW: FORTRAN90, C, MPI
- •ACRES: FORTRAN90, C, HPF





Performance

- Beta Testing Complete
- Scalability: (32 P.E.)

O2K	31	25	
SP	52	47	
	DoDPW	ACRES	

Management

(top view)

- Leveraging: NRL, ONR (Harvard)
- Transition Approach: Code supported by
- Funding \$1,672K (FY96-00)

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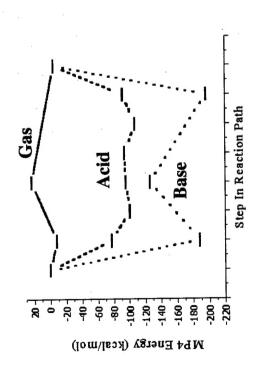
Computational Chemistry and Materials Science

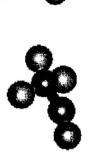
CCM-2: Quantum Chemistry

Synopsis

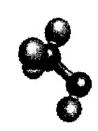
- codes to investigate new chemical species Objective: Produce ab initio chemistry
- Partners: AFRL/ML, U. of Delaware, Iowa State U.
- Development Paradigm: FORTRAN77,

Reaction Pathways - Formation of POSS









Base Catalysis - POSS synthesis

Performance

- Beta Testing Complete
- 57% (SP) 50% (O2K) Scalability:
- Operating Platforms: SP, O2K

Management

- Leveraging: DoE, NSF
- Transition Approach: Used by 1,000+ users world-wide
- Funding \$1,593K (FY96-00)

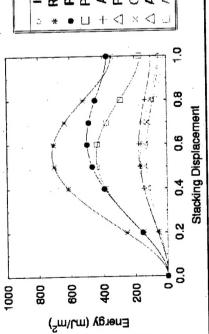
olimits CM Computational Chemistry and Materials Science

CCM-3: Tight Binding Molecular Dynamics

- Synopsis
- *Objective*: Develop a general purpose tight-binding MD code and associated libraries
- Partners: Ohio State University, George Mason University to be consistent of others.
 - Development Paradigm: FORTRAN77, MPI

First Principles Calculation of Mechanical Properties: Stacking Fault Energies and Ductility Criterion

D. A. Papaconstantopoulos and M. J. Mehl Center for Computational Materials Science Naval Research Laboratory Washington DC



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- Performance
- Beta Testing Complete
- Scalability: (16 P.E.)

	SP	02K
Fitting	1	40
Static	50	95
TBMD	62	76

Management

- Leveraging: NRL
- Transition Approach: Supported by NRL
- Funding \$1,954K (FY96-00)

CCM

Computational Chemistry and Materials Science

CCM-4: Classical Molecular Dynamics

Synopsis

- Objective: Develop a molecular dynamics package that can realistically simulate large molecular systems for long time scales
- Partners: AFRL, ARL, Iowa State U_GU. of Utah, U_Sof Michigan, U. of Houston, Cornell U.
 - Development Paradigm: FORTRAK77, MPI

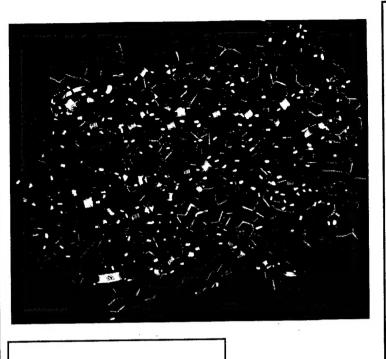
Performance

- Beta Testing Complete
- Scalability:

	FMD	СМД	
T3E	40	1	
SP	40	85	
O2K	44	80	

Management

- Leveraging: AFRL, AFOSR
- Transition Approach: Supported by AFRL
- Funding \$1,793K (FY96-00)



Applications of CHSSI Technology

Problem

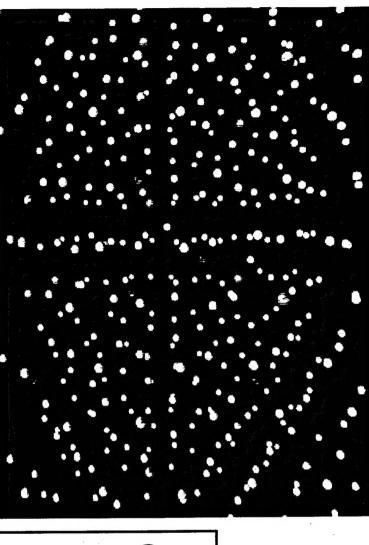
- Objective: Is there global recombination of boron atoms in a solid hydrogen matrix if a single recombination is initiated?
- Technology Used: CMD (CCM-4)
 - Technical Team: U. of Utah,

AFRL/PR

Hydrogen O art mas sufferent?

Non-reacting Boron atom O

Reacting Boron atom O



■ Impact...

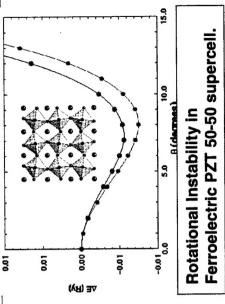
- of Code Application: The thermal conductivity of the hydrogen matrix is sufficient to prevent global recombination of boron atoms in solid hydrogen
 - of CHSSI: This calculation would be take over *X* year, possibly two, to complete. The parallel CMD performs the calculations in β' weeks on 64 nodes of an IBM SP.

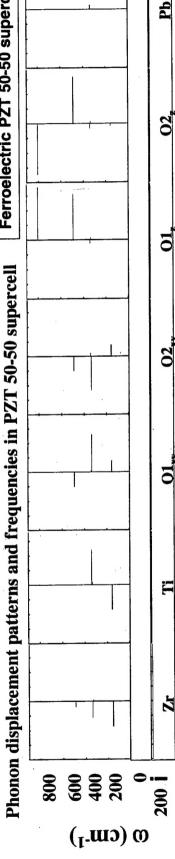
CM

Applications of CHSSI Technology



- Objective: Determine if Pb can be removed from PZT [Pb(Zr,Ti)O₃] a Navy Sonar Material, and maintain its effectiveness
- Technology Used: DoDPW, NRL-LAPW
- Technical Team: NRL, ONR





Impact...

- of Code Application: ONR and NRL are gaining a crucial understanding of how the Pb effects the sonar material
- of CHSSI: The code accurately describes the material and researchers will be able to suggest effective substitutes within a short time

CM

Applications of CHSSI Technology

Problem

- *Objective*: Identify important factors in the synthesis of polyhedral oligomeric silsesquioxanes (POSS) to modify the existing lengthy synthesis process
 - Technology Used: GAMESS (CCM-2)
- Technical Team: AFRL, Iowa State U., ASC MSRC

Impact...

- of Code Application: The effects of solvents led to a new process for the formulation of POSS. of different catalysts and solvents have been found to be crucial in the quick formulation of POSS.
- of CHSSI: The calculation to determine the structure and associated energies lasted approximately $\mathcal{X}^{\text{week}}$. The serial version would require over ONE YEAR of dedicated time.

